

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081721\
 Data File : BM031795.D
 Acq On : 17 Aug 2021 23:31
 Operator : CG/JU
 Sample : M3425-12
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 22N

Quant Time: Aug 18 02:26:05 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 16 16:42:38 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.545	152	28777	20.000	ng	# 0.00
21) Naphthalene-d8	10.304	136	113170	20.000	ng	#-0.01
39) Acenaphthene-d10	14.180	164	74167	20.000	ng	-0.01
64) Phenanthrene-d10	16.939	188	162311	20.000	ng	0.00
76) Chrysene-d12	21.133	240	155304	20.000	ng	-0.01
86) Perylene-d12	23.285	264	143653	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.187	112	278958	155.585	ng	0.00
7) Phenol-d6	6.745	99	309599	119.166	ng	0.00
23) Nitrobenzene-d5	8.698	82	268119	99.200	ng	0.00
42) 2,4,6-Tribromophenol	15.686	330	134495	147.236	ng	0.00
45) 2-Fluorobiphenyl	12.798	172	540940	97.078	ng	0.00
79) Terphenyl-d14	19.586	244	1014103	110.018	ng	0.00

Target Compounds	Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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